

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042423\
 Data File : BG057158.D
 Acq On : 24 Apr 2023 19:32
 Operator : CG/JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC060

Manual Integrations

APPROVED

Reviewed By : Christian Giraldo 04/25/2023
 Supervised By : Jagrut Upadhyay 04/25/2023

Quant Time: Apr 25 04:22:28 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG042423.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Tue Apr 25 04:18:38 2023

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.390	152	27027	20.000	ng	0.00
21) Naphthalene-d8	11.227	136	108258	20.000	ng	0.00
39) Acenaphthene-d10	15.004	164	77446	20.000	ng	0.00
64) Phenanthrene-d10	17.742	188	180057	20.000	ng	0.00
76) Chrysene-d12	22.060	240	157437	20.000	ng	0.00
86) Perylene-d12	25.584	264	160742	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.899	112	191053	128.102	ng	0.00
7) Phenol-d6	7.526	99	282820	131.064	ng	0.00
23) Nitrobenzene-d5	9.571	82	294010	134.261	ng	0.00
42) 2,4,6-Tribromophenol	16.485	330	112653	110.246	ng	0.00
45) 2-Fluorobiphenyl	13.630	172	658916	114.341	ng	0.00
79) Terphenyl-d14	20.315	244	1064344	118.248	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.708	88	43408	64.952	ng	94
3) Pyridine	4.131	79	141760	69.911	ng	95
6) Aniline	7.702	93	184113	65.334	ng	97
8) 2-Chlorophenol	7.949	128	97641	62.495	ng	98
9) Benzaldehyde	7.509	77	75459	58.383	ng	87
10) Phenol	7.550	94	139812	65.252	ng	99
11) bis(2-Chloroethyl)ether	7.791	93	110346	65.628	ng	92
12) 1,3-Dichlorobenzene	8.278	146	106215	57.388	ng	96
13) 1,4-Dichlorobenzene	8.425	146	107403	57.581	ng	99
14) 1,2-Dichlorobenzene	8.748	146	103193	57.300	ng	96
15) Benzyl Alcohol	8.619	79	109888	69.110	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.907	45	144322m	63.285	ng	
17) 2-Methylphenol	8.824	107	102174	65.165	ng	97
18) Hexachloroethane	9.488	117	39316	62.132	ng	90
20) 3+4-Methylphenols	9.153	107	138452	64.524	ng	99
22) Acetophenone	9.218	105	183053	62.331	ng	# 94
24) Nitrobenzene	9.612	77	150805	68.590	ng	99
25) Isophorone	10.134	82	285155	67.889	ng	98
26) 2-Nitrophenol	10.328	139	52280	63.647	ng	90
27) 2,4-Dimethylphenol	10.369	122	99953	63.313	ng	98
28) bis(2-Chloroethoxy)met...	10.604	93	146855	66.696	ng	97
29) 2,4-Dichlorophenol	10.863	162	106461	60.535	ng	96
30) 1,2,4-Trichlorobenzene	11.086	180	118225	54.974	ng	95
31) Naphthalene	11.280	128	341425	59.378	ng	99
32) Benzoic acid	10.522	122	46271	60.625	ng	92
33) 4-Chloroaniline	11.380	127	152858	60.463	ng	99
34) Hexachlorobutadiene	11.550	225	82658	52.125	ng	97
35) Caprolactam	12.155	113	32660	67.067	ng	91
36) 4-Chloro-3-methylphenol	12.472	107	120597	63.743	ng	95
37) 2-Methylnaphthalene	12.854	142	265114	58.590	ng	96
38) 1-Methylnaphthalene	13.072	142	247359	59.211	ng	98
40) 1,2,4,5-Tetrachloroben...	13.213	216	158130	55.801	ng	98
41) Hexachlorocyclopentadiene	13.189	237	69720	58.849	ng	96
43) 2,4,6-Trichlorophenol	13.448	196	94711	61.734	ng	95
44) 2,4,5-Trichlorophenol	13.524	196	112082	60.514	ng	95
46) 1,1'-Biphenyl	13.841	154	342596	59.348	ng	97

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042423\
 Data File : BG057158.D
 Acq On : 24 Apr 2023 19:32
 Operator : CG/JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC060

Manual Integrations

APPROVED

Reviewed By :Christian Giraldo 04/25/2023
 Supervised By :Jagrut Upadhyay 04/25/2023

Quant Time: Apr 25 04:22:28 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG042423.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Tue Apr 25 04:18:38 2023

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
47) 2-Chloronaphthalene	13.894	162	258449	59.319	ng	98
49) Acenaphthylene	14.734	152	418520	60.511	ng	99
50) Dimethylphthalate	14.440	163	333326	61.326	ng	100
51) 2,6-Dinitrotoluene	14.570	165	74864	65.986	ng	98
52) Acenaphthene	15.069	154	275752m	61.020	ng	
53) 3-Nitroaniline	14.904	138	73746	65.458	ng	95
54) 2,4-Dinitrophenol	15.110	184	37707	60.360	ng	94
55) Dibenzofuran	15.398	168	422655	57.725	ng	97
56) 4-Nitrophenol	15.198	139	52875	66.678	ng	90
57) 2,4-Dinitrotoluene	15.351	165	102035	66.039	ng	96
58) Fluorene	16.044	166	338880	57.487	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.621	232	94986m	59.894	ng	
60) Diethylphthalate	15.786	149	326342	62.287	ng	99
61) 4-Chlorophenyl-phenyle...	16.021	204	196932	55.660	ng	96
62) 4-Nitroaniline	16.062	138	73928	65.168	ng	92
63) Azobenzene	16.314	77	372880	66.903	ng	97
65) 4,6-Dinitro-2-methylph...	16.120	198	58271	60.895	ng	97
66) n-Nitrosodiphenylamine	16.238	169	292811	62.324	ng	99
67) 4-Bromophenyl-phenylether	16.919	248	124715	57.596	ng	96
68) Hexachlorobenzene	17.049	284	121711	54.801	ng	96
69) Atrazine	17.166	200	94380	62.895	ng	95
70) Pentachlorophenol	17.389	266	68541	65.472	ng	98
71) Phenanthrene	17.789	178	545848	59.383	ng	100
72) Anthracene	17.877	178	553851	59.590	ng	99
73) Carbazole	18.141	167	478456	60.985	ng	99
74) Di-n-butylphthalate	18.658	149	477696	65.457	ng	99
75) Fluoranthene	19.774	202	663952	58.139	ng	99
77) Benzidine	19.939	184	140998	59.433	ng	99
78) Pyrene	20.139	202	668332	62.525	ng	99
81) Benzo(a)anthracene	22.042	228	653876	60.406	ng	99
82) 3,3'-Dichlorobenzidine	21.936	252	181716	64.140	ng	96
83) Chrysene	22.112	228	618964	59.294	ng	99
84) Bis(2-ethylhexyl)phtha...	21.889	149	272315	60.724	ng	98
85) Di-n-octyl phthalate	23.193	149	470468	66.628	ng	97
87) Indeno(1,2,3-cd)pyrene	29.661	276	704684	61.201	ng	# 95
88) Benzo(b)fluoranthene	24.456	252	619665	61.498	ng	100
89) Benzo(k)fluoranthene	24.533	252	648498	62.220	ng	98
90) Benzo(a)pyrene	25.420	252	614061	61.874	ng	99
91) Dibenzo(a,h)anthracene	29.726	278	580790	60.749	ng	97
92) Benzo(g,h,i)perylene	30.953	276	570971	60.361	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

